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WALPOLE ISLAND WATER TREATMENT PLANT

ANNUAL REPORT 1990

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DRINKING WATER SURVEILLANCE PROGRAM

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JULY 1992



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EXECUTIVE SUMMARY

DRINKING WATER SURVEILLANCE PROGRAM

WALPOLE ISLAND WATER TREATMENT PLANT 1990 ANNUAL REPORT

The Drinking Water Surveillance Program (DWSP) for Ontario is a monitoring program providing immediate, reliable, current information on drinking water quality. The DWSP officially began in April 1986 and is designed to eventually include all municipal supplies in Ontario. In 1990, 76 systems were being monitored.

The Walpole Island water treatment plant is a package plant which uses conventional treatment and treats water from the St. Clair River. The process consists of coagulation, flocculation, sedimentation, filtration, taste and odour control and disinfection. This plant has a rated capacity of 0.9 x 1000 m^3/day . The Walpole Island water treatment plant serves a population of approximately 2,900.

Raw and treated water at the plant was sampled for the presence of approximately 180 parameters. Parameters were divided into the following groups: bacteriological, inorganic and physical (laboratory chemistry, field chemistry and metals), and organic (chloroaromatics, chlorophenols, pesticides and PCB, phenolics, polyaromatic hydrocarbons, specific pesticides and volatiles). Samples were analyzed for specific pesticides and chlorophenols twice a year in the spring and fall.

Table A is a summary of all results by group.

No known health related guidelines were exceeded.

The Walpole Island water treatment plant, for the sample year 1990, produced good quality water. Water from the distribution system was not sampled.

TABLE A DRINKING WATER SURVEILLANCE PROGRAM WALPOLE ISLAND WTP

SUMMARY TABLE BY SCAN

A POSITIVE VALUE DEMOTES THAT THE RESULT IS CREATER THAN THE STATISTICAL LIMIT OF DETECTION AND IS QUANTIFIABLE A '.' INDICATES THAT NO SAMPLE WAS TAKEN SITE SITE

	SCAN	TESTS	POSITIVE	RAU %POSITIVE	TESTS	POSITIVE %	TED POSITIVE	
	BACTERIOLOGICAL	m	0	0	2	-	50	
	CHEMISTRY (FLD)	18	18	100	36	36	100	
	CHEMISTRY (LAB)	132	8	75	131	92	0.2	
	METALS	144	36	ر 25	144	39	27	
	CHLOROAROMATICS	8	0	0	78	0	0	
	CHLOROPHENOLS	12	0	0	12	0	0	
	РАН	102	0	0	102	0	0	
	PESTICIDES & PCB	204	0	0	204	0	0	
	PHENOL 1 CS	9	0	0	9	0	0	
	SPECIFIC PESTICIDES	58	0	0	50	0	0	
	VOLATILES	174	0	0	145	20	13	
TOTAL		937	153		916	188		

TOTAL

WALPOLE ISLAND WATER TREATMENT PLANT 1990 ANNUAL REPORT

INTRODUCTION

The Drinking Water Surveillance Program (DWSP) for Ontario is a monitoring program providing immediate, reliable, current information on drinking water quality. The DWSP officially began in April 1986 and is designed to eventually include all municipal supplies in Ontario. In 1990, 76 systems were being monitored.

Appendix A has a full description of the DWSP.

The DWSP was initiated for the Walpole Island water treatment plant in the spring of 1985 as part of a survey of the St.Clair /Detroit River area. Previous DWSP annual reports have been published for 1986, 1987, 1988 and 1989.

PLANT DESCRIPTION

The Walpole Island water treatment plant is a package plant which uses conventional treatment and treats water from St. Clair River. The process consists of coagulation, flocculation, sedimentation, filtration, taste and odour control, and disinfection. This plant has a rated capacity of 0.9 x 1000 m^3/day . The Walpole Island water treatment plant serves a population of approximately 2,900.

The sample day flows ranged from 0.4 x 1000 m^3/day to 0.5 x 1000 m^3/day .

General plant information is presented in Table 1 and a schematic of plant processes, chemical addition points and sampling locations in Figure 1.

SAMPLING AND ANALYSES

Sample lines in the plant were flushed prior to sampling to ensure that the water obtained was indicative of its origin and not residual water standing in the sample line.

At all distribution system locations two types of samples were obtained, a standing and a free flow. The standing sample consisted of water that had been in the household plumbing and service connection for a minimum of six hours. These samples were used to make an assessment of the change in the levels of inorganic compounds and metals, due to leaching from, or deposition on, the plumbing system. The only analyses carried out on the standing samples therefore, were General Chemistry and Metals. The free flow sample represented fresh water from the distribution main, since the sample tap was flushed for five minutes prior to sampling.

Attempts were made to capture the same block of water at each sampling point by taking the retention time into consideration. Retention time was calculated by dividing the volume of water between two sampling points by sample day flow. For example, if it was determined that retention time within the plant was five hours, then there would be a five hour interval between the raw and treated sampling. Similarly, if it was estimated that it took approximately one day for the water to travel from the plant to the distribution system site, this site would be sampled one day after the treated water from the plant.

Stringent DWSP sampling protocols were followed to ensure that all samples were taken in a uniform manner (see Appendix B).

Plant operating personnel routinely analyze parameters for process control (Table 2).

Raw and treated water at the plant was sampled for the presence of approximately 180 parameters. Parameters were divided into the following groups: bacteriological, inorganic and physical (laboratory chemistry, field chemistry and metals), and organic (chloroaromatics, chlorophenols, pesticides and PCB, phenolics, polyaromatic hydrocarbons, specific pesticides and volatiles). Samples were analyzed for specific pesticides and chlorophenols twice a year in the spring and fall. Laboratory analyses were conducted at the Ministry of the Environment facilities in Rexdale, Ontario.

RESULTS

Field measurements were recorded on the day of sampling and were entered onto the DWSP database as submitted by plant personnel.

Table 3 contains information on delay time between raw and treated water sampling, flow rate, and treatment chemical dosages.

Table 4 is a summary break-down of the number of water samples analyzed by parameter and by water type. The number of times that a positive or trace result was detected is also reported.

Positive denotes that the result is greater than the statistical limit of detection established by the Ministry of the Environment laboratory staff and is quantifiable. Trace (<T) denotes that the level measured is greater than the lowest value detectable by the method but lies so close to the detection limit that it cannot be confidently quantified.

Table 5 presents the results for parameters detected on at least one occasion.

Table 6 lists all parameters analyzed in the DWSP.

Associated guidelines and detection limits are also supplied on Tables 5 and 6. Parameters are listed alphabetically within each scan.

DISCUSSION

GENERAL

Water quality was judged by comparison with the Ontario Drinking Water Objectives publication (ODWOs). When an Ontario Drinking Water Objective (ODWO) was not available, guidelines/limits from other agencies were used. These guidelines were obtained from the Parameter Listing System database.

IN THIS REPORT, DISCUSSION IS LIMITED TO:

- THE TREATED AND DISTRIBUTED WATER;
- ONLY THOSE PARAMETERS WITH CONCENTRATIONS ABOVE GUIDELINE VALUES; AND
- POSITIVE ORGANIC PARAMETERS DETECTED.

BACTERIOLOGICAL

Guidelines for bacteriological sampling and testing of a supply are developed to maintain a proper supervision of its bacteriological quality. Routine monitoring programs usually require that multiple samples be collected in a given system. Full interpretation of bacteriological quality cannot be made on the basis of single samples.

Standard plate count was the only bacteriological analysis conducted on the treated samples. No results were reported above the guideline.

INORGANIC & PHYSICAL

CHEMISTRY (FLD)

It is desirable that the temperature of drinking water be less than 15° C. The palatability of water is enhanced by its coolness. A temperature below 15° C will tend to reduce the growth of nuisance organisms and hence minimize associated taste, colour, odour and corrosion problems. The temperature of the delivered water may increase in the distribution system due to the warming effect of

the soil in late summer and fall and/or as a result of higher temperatures in the source water.

Field temperature exceeded the ODWO Maximum Desirable Concentration of 15° C in 2 of 6 treated water samples with a maximum reported value of 21.2° C.

CHEMISTRY (LAB)

The ODWOs indicate that a hardness level of between 80 and 100 mg/L as calcium carbonate for domestic waters provides an acceptable balance between corrosion and encrustation. Water supplies with a hardness greater than 200 mg/L are considered poor and would possess a tendency to form scale deposits and result in excessive soap consumption.

Hardness exceeded the ODWO Aesthetic or Recommended Operational Guideline of 80-100 mg/L in the 6 treated water samples with a maximum reported value of 121.9 mg/L.

METALS

At present, there is no evidence that aluminum is physiologically harmful and no health limit for drinking water has been specified. The measure of aluminum in treated water is important to indicate the efficiency of the treatment process. The ODWOs indicate that a useful guideline is to maintain a residual below 100 ug/L as aluminum in the water leaving the plant, to avoid problems in the distribution system.

Aluminum exceeded the ODWO Aesthetic or Recommended Operational Guideline of 100 ug/L in 3 of 6 treated water samples with a maximum reported value of 270.0 ug/L.

ORGANIC

CHLOROAROMATICS

The results of the chloroaromatic scan showed that none were detected.

CHLOROPHENOLS

The results of the chlorophenol showed that none were detected.

POLYAROMATIC HYDROCARBONS (PAH)

The results of the PAH scan showed that none were detected.

PESTICIDES & PCB

The results of the PCB scan showed that none were detected.

The results of the regular pesticide scan showed that none were detected above trace levels.

PHENOLICS

Phenolic compounds are present in the aquatic environment as a result of natural and/or industrial processes. The ODWOS recommend, as an operational guideline, that phenolic substances in drinking water not exceed 2.0 ug/L. This limit has been set primarily to prevent undesirable taste and odours, particularly in chlorinated water. No results were reported above trace levels.

SPECIFIC PESTICIDES

The results of the specific pesticides scan showed that none were detected.

VOLATILES

The detection of benzene, ethylbenzene, toluene and xylenes at low, trace levels may be a laboratory artifact derived from the analytical methodology.

Trihalomethanes (THMs) are produced during the water treatment process and will always occur in chlorinated waters. THMs are comprised of chloroform, chlorodibromomethane and dichlorobromomethane; bromoform occurs occasionally. Results are reported for the individual compounds as well as for total THMs. Only total THMs results are discussed.

Total THMs were found at positive levels in the 5 treated water samples analyzed with a maximum level of 32.6 ug/L. This was below the ODWO Maximum Acceptable Concentration of 350 ug/L.

CONCLUSIONS

The Walpole Island water treatment plant, for the sample year 1990, produced good quality water. Water from the distribution system was not sampled.

No known health related guidelines were exceeded.

FIGURE 1

WALPOLE ISLAND WATER TREATMENT PLANT



TABLE 1

DRINKING WATER SURVEILLANCE PROGRAM

PLANT GENERAL REPORT

WORKS #: 230000129 PLANT NAME: WALPOLE ISLAND

DISTRICT: WALPOLE ISLAND REGION: SOUTHWESTERN DISTRICT OFFICER: O. WIGLE

UTM #:

PLANT SUPERINTENDENT: J. TOOSHKENING

ADDRESS: WALPOLE ISLAND, ONTARIO N8A 4K9

(Telephone) (519 627-1426)

MUNICIPALITY: WALPOLE ISLAND AUTHORITY: FEDERAL GOVERNMENT

PLANT INFORMATION:

PLANT VOLUME: 0.000	(X	1000 m3	3)		
DESIGN CAPACITY:		2.500	(X	1000	m3/day)
RATED CAPACITY:		0.000	(X	1000	m3/day)

MUNICIPALITY: WALPOLE ISLAND (SUMMER MONTHS) **POPULATION:** 1,900 2,100

TABLE 2 DRINKING WATER SURVEILLANCE PROGRAM IN-PLANT MONITORING

PARAMETER	LOCATION	FREEDENCY
FREE CHLORINE RESIDUAL	TREATED WATER IN LAB RAW WATER IN LAB	2/DAY 2/DAY
TOTAL CHLORINE RESIDUAL	TREATED WATER IN LAB SETTLED WATER IN LAB	2/DAY 2/DAY
PH	TREATED WATER IN LAB RAW WATER IN LAB	WEEKLY WEEKLY
TURBIDITY	TREATED WATER IN LAB SETTLED WATER IN LAB RAW WATER IN LAB	2/DAY 2/DAY 2/DAY

TABLE 3 DRINKING WATER SURVEILLANCE PROGRAM WALPOLE ISLAND WTP SAMPLE DAY CONDITIONS FOR 1990

POST CHLORINATION	CHLORINE	1.00 .50 .50 .50 .50
PRE CHLORINATION	CHLORINE	. 80 . 80 . 50 . 80 . 70
COAGULATION AID	POLYELECTROLYTE	25 25 10 10 10
ILCAL DOSAGES (MG/L) TASTE & ODOUR	ACTIVATED CARBON POWDER	10.00 10.00 10.00 10.00 10.00
TREATMENT CHEN COAGULATION	ALUM DRY	7.50 10.00 7.50 7.50 7.50
	FLOW (1000M3)	.408 .534 .475
	DELAY * TIME(HRS)	3 48.73 5 49.50 7 51.51 4 .37 6 41.86 5 20.00
	DATE	MAR 0 MAY 0 SEP 0

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* THE DELAY TIME BETWEEN THE RAW AND TREATED WATER SAMPLING, SHOULD ESTIMATE THE RETENTION TIME.

Page 9

		1	RAW		TREA	FED
SCAN PARAMETER	TOTAL POS	ITIVE TR	ACE	TOTAL POS	ITIVE TR	ACE
BACTERIOLOGICAL						
FECAL COLIFORM MF	1	0	0	:	:	:
STANDRD PLATE CNT MF	;	÷	÷	2	1	0
T COLIFORM BCKGRD MF	i	D	0		:	
*TOTAL GROUP BACTERIO	LOGICAL 3	0	0	2	1	0
CHEMISTRY (FLD)						
FLD CHLORINE (COMB)	•	•	•	6	6	0
FLD CHLORINE FREE	•		:	6	6	õ
FLD PH	6	6	0	6	6	0
FLD TEMPERATURE	6	6	0	6	6	0
FLD TURBIDITY	6	6	0	6	0	U
*TOTAL SCAN CHEMISTRY	(FLD) 18	18	0	36	36	0
CHEMISTRY (LAB) ALKALINITY CALCIUM CYANIDE COLOUR COLOUR CONDUCTIVITY DISS ORG CARBON FLUORIDE HARDNESS IONCAL LANGELIERS INDEX MAGNESIUM SOOIUM AMMONIUM TOTAL NITRITE TOTAL NITRATES NITROGEN TOT KJELD PH	6 6 6 6 6 6 6 6 6 6 6 6 6 6 6 6	6 6 0 6 6 6 6 6 6 6 6 6 6 6 6 6 6 6 6 6	0005000000260003	6 6 6 6 6 6 6 6 6 6 6 6 6 6 6 6 6 6 6	6 6 6 6 6 6 6 6 6 6 6 0 0 6 4 6 1	00030000000140202
PHOSPHORUS FIL REACT PHOSPHORUS TOTAL SULPHATE	6	0	6	6	0	307
TURBIDITY	6	6	0	0	3	2
*TOTAL SCAN CHEMISTR	Y (LAB) 132	99	22	131	92	18

			RAW		TR	EATED
SCAN		0001711/5			0001711/5	70405
PARAMETER	TOTAL	PUSITIVE	TRACE	TOTAL	PUSITIVE	TRACE
METALS						
SILVER	6	0	0	6	0	0
ALUMINUM	6	6	0	6	6	0
ARSENIC	6	0	6	6	0	6
BARIUM	6	6	0	6	6	D
BORON	0	2	4	0	1	2
CADMILIN	6	ň	1	6	n N	1
COBALT	6	ŏ	6	6	õ	Ś
CHROMIUM	6	Ō	4	6	Ō	4
COPPER	6	D	6	6	1	5
IRON	6	1	5	6	D	1
MERCURY	6	0	1	6	0	1
MANGANESE	6	6	0	6	5	1
NICKEL	6	3	3	6	4	2
LEAD	6	0	6	6		6
ANTIMONY	6	ĩ	5	6	1	5
SELENIUM	6	ó	2	6	1	1
STRONTIUM	6	6	D	6	6	0
TITANIUM	6	3	3	6	1	5
THALLIUM	6	0	0	6	0	0
URANIUM	6	0	6	6	0	5
ZINC	6	0	6	6	0	6
ZINC	0	۷	4	0	0	U
*TOTAL SCAN METALS						
	144	36	71	144	39	61
*TOTAL GROUP INORGANIC	: & PH'	SICAL				
	294	153	93	311	167	79
CHLOROAROMATICS						
HEXACHLOROBUTADIENE	6	. 0	0	6	0	0
123 TRICHLOROBENZENE	6	0	0	6	0	0
1234 T-CHLOROBENZENE	6	0	0	6	0	0
1235 T-CHLOROBENZENE	6	0	0	6	0	0
124 TRICHLOROBENZENE	0 4	0	0	0 4	U	0
135 TRICHLOROBENZENE	6	ő	0	6	0	ñ
HCB	6	ŏ	ŏ	6	Ő	ŏ
HEXACHLOROETHANE	6	0	D	6	0	0
OCTACHLOROSTYRENE	6	0	D	6	0	0
PENTACHLOROBENZENE	6	0	0	6	0	0
236 TRICHLOROTOLUENE	6	0	0	6	0	0
245 TRICHLOROTOLUENE	6	0	0	6	0	0
ZOW TRICHLORUTULUENE	6	U	0	0	U	U
TOTAL SCAN CHLOROARON	ATICS					
COLLE COM CHECKOAKO	84	0	0	84	0	0

CHLOROPHENOLS

.

			RAW		TREAT	ED
SCAN	70741 00011		ACE 1		TIVE TO	CE
PARAMETER	TUTAL POST	IVE IN	ALE	IUTAL POS		
234 TRICHLOROPHENOL	2	0	0	2	0	0
2345 T-CHLOROPHENOL	2	0	0	2	0	0
2356 T-CHLOROPHENOL	2	0	0	2	0	0
245-TRICHLOROPHENOL	2	0	0	2	0	0
246-TRICHLOROPHENDL	2	0	0	2	0	0
PENTACHLOROPHENOL	۷	U	0	2	v	0
*TOTAL SCAN CHLOROPHE	NOLS					
	12	0	0	12	0	0
РАН						
				,	•	•
PHENANTHRENE	6	0	0	0	0	0
ANTHRACENE	o 4	0	0	6	ñ	ň
PLOUKANTHENE	6	0	ñ	6	ñ	ŏ
RENZOLAJANTHRACENE	6	ŏ	ŏ	6	ŏ	ŏ
CHRYSENE	6	Ō	Ō	6	0	0
DIMETH. BENZ(A)ANTHR	6	Ō	Ó	6	0	0
BENZO(E) PYRENE	6	0	0	6	0	0
BENZO(B) FLUORANTHEN	6	0	0	6	0	0
PERYLENE	6	0	0	6	0	0
BENZO(K) FLUORANTHEN	6	0	0	6	0	0
BENZO(A) PYRENE	6	0	0	6	0	0
BENZO(G,H,I) PERYLEN	°,	0	0	٥ 4	0	0
DIBENZO(A,H) ANTHRAC	0 4	0	0	6	ň	ň
PENZO(P) CHOYSENE	6	n	ň	6	ő	ŏ
CORONENE	6	ŏ	ŏ	6	ō	Õ
*TOTAL SCAN PAH	402	•	•	103	0	0
	102	U	U	102	U	0
PESTICIDES & PCB						
ALDRIN	6	0	0	6	0	0
ALPHA BHC	6	0	5	6	0	5
BETA BHC	6	0	0	6	0	0
LINDANE	6	0	0	6	0	0
ALPHA CHLORDANE	6	0	0	°,	U	0
GAMMA CHLORDANE	°,	0	0	0 4	0	0
DIELDKIN	6	0	ň	6	ñ	ň
ENDOSIU EAN 1	6	ñ	ň	6	ő	ŏ
	6	ŏ	ŏ	6	õ	ŏ
ENDRIN	6	õ	Ő	6	0	0
ENDOSULFAN SULPHATE	6	0	0	6	0	0
HEPTACHLOR EPOXIDE	6	0	0	6	0	0
HEPTACHLOR	6	0	0	6	0	0
MIREX	6	0	0	6	0	0
OXYCHLORDANE	6	0	0	6	0	0
OPDDT	6	0	0	6	0	0
PCB	0	0	0	6	0	0
PPDDE	6	0	0	6	õ	ŏ
1 TODE	0	~			•	-

			RAW		TF	REATED
SCAN PARAMETER	TOTAL	POSITIVE	TRACE	TOTAL	POSITIVE	TRACE
PPDDT	6	0	0	6	0	0
AMETRINE	6	0	0	6	0	0
ATRAZINE	6	0	0	6	0	0
ATRATONE	6	0	0	6	0	0
CYANAZINE (BLADEX)	6	0	0	6	0	0
DESETHYLATRAZINE	0	0	U	0	0	0
DPONETONE	2	0	0	5	0	0
PROPAZINE	6	0	0	6	0	ő
PROMETRYNE	6	ŏ	ŏ	6	ŏ	ŏ
METRIBUZIN (SENCOR)	6	Ō	0	6	0	0
SIMAZINE	6	0	0	6	0	0
ALACHLOR (LASSO)	6	0	0	6	0	0
METOLACHLOR	6	0	0	6	0	0
HEXACLCYCLOPENTADIEN	1	0	0	1	0	0
TOTAL SCAN DESTICIDE	s 2 pri	2				
TOTAL SCAN PESTICIDE	204	, 0	5	204	0	5
			-			-
PHENOLICS						
PHENOLICS	6	0	1	6	0	1
*TOTAL SCAN PHENOLICS						
	6	0	1	6	0	1
SPECIFIC PESTICIDES						
TOXAPHENE	6	0	0	6	0	0
2,4,5-T	2	0	0	2	0	0
2,4-0	2	0	0	2	0	0
2,4-DB	2	0	0	2	0	0
DICAMBA	2	0	0	2	0	ñ
PICHLORAM	0	ő	ő	0	ŏ	ŏ
SILVEX	2	ŏ	ŏ	2	ŏ	ŏ
DIAZINON	2	Ō	Ō	2	Ő	Ō
DICHLOROVOS	2	0	0	2	0	0
CHLORPYRIFOS	2	0	0	2	0	0
ETHION	2	0	0	2	0	0
AZINPHOS-METHYL	0	0	0	0	0	0
MALATHION	2	0	0	2	0	U
MEVINPHOS	2	U	0	2	0	0
METHYL PAKATHION	2	0	0	2	0	0
PARATHION	2	0	0	2	0	ő
PHORATE	2	0	0	2	ő	ŏ
RELDAN	2	Ő	Ő	2	Ő	D
RONNEL	2	Ő	Ő	2	0	0
AMINOCARB	0	0	0	0	0	0
BENONYL	0	0	0	0	0	0
BUX	0	0	0	0	0	0
CARBOFURAN	2	0	0	1	0	0
CICP	2	0	0	1	0	0
DIALLAIE	2	0	U	1	U	v

Ψ.

				RAW		1	IREATE	D
SCAN PAPAMETER	TOTAL	POSITIN	/E	TRACE	TOTAL	POSITIVE	TRAC	E
EPTAM	2		0	0	1	(2	0
IPC	2		0	0	1		5	0
PROPOXUR	2		0	0	1	L. L.	J	0
CARBARYL	2		C	0	1		J	0
BUTYLATE	2		0	0	1	, i	J	υ
*TOTAL SCAN SPECIFIC	PESTIC	IDES						
	58		0	0	50	(D	0
VOLATILES								
BENZENE	6		0	4	5	(0	4
TOLUENE	6		0	3	5		0	4
ETHYLBENZENE	6		0	2	5		0	4
P-XYLENE	6		0	0	5		0	0
M-XYLENE	6		0	1	5		D	2
O-XYLENE	6		0	1	5		0	1
STYRENE	6	•	0	4	2		U	U
1,1 DICHLOROETHYLENE	6		0	0	2		U	0
METHYLENE CHLORIDE	6	1	0	1	2		U	U
T1, 2DICHLOROETHYLENE	6	•	0	0	2		0	0
1,1 DICHLOROETHANE	6)	0	U	2		5	0
CHLOROFORM	6	2	0	1	2		2	1
111, TRICHLOROETHANE	ć	•	0	1	2		0	
1,2 DICHLOROETHANE	6)	0	0	2		0	0
CARBON TETRACHLORIDE	9)	U	0	2		0	0
1,2 DICHLOROPROPANE	9)	U	0	2		0	0
TRICHLOROETHYLENE	9	2	0	U	2		5	0
DICHLOROBROMOMETHANE	2)	0	0	2		5	0
112 TRICHLOROETHANE)	U	0	2		5	0
CHEDRODIBROMOMETHANE	2		0	1	2		0	0
I-CHLOROETHTLENE	2		0		2		0	5
BRUMUFURM) ,	0		2		0	5
1122 1-CHLOROETHANE	2		0	0	2		0	ň
CHLOROBENZENE	2)	0	0	2		0	0
1,4 DICHLOROBENZENE			0	0	5		ñ	ñ
1,5 DICHLOROBENZENE		2	0	0			ň	ň
1,2 DICHLOROBENZENE		, ,	0	0	2		0	ñ
ETHLITENE DIBROMIDE		2	0	0	2		5	ñ
TOTE TRINALOMETHANES	c	>	U	U	2		2	0
*TOTAL SCAN VOLATILES	5			10	4/5	-	20	21
	174	•	0	19	143		.0	- 1
TUTAL GROUP ORGANIC	641)	0	25	603	; 2	20	27

KEY TO TABLE 5 and 6

- ONTARIO DRINKING WATER OBJECTIVES (ODWO) A
 - 1. Maximum Acceptable Concentration (MAC)
 - 1+. MAC for Total Trihalomethanes
 - 2. Interim Maximum Acceptable Concentration (IMAC)

 - Aesthetic Objective (AO)
 AO for Total Xylenes
 Recommended Operational Guideline
- HEALTH & WELFARE CANADA (H&W) B
 - Maximum Acceptable Concentration (MAC)
 Proposed MAC
 Interim MAC

 - 4. Aesthetic Objective (AO)
- WORLD HEALTH ORGANIZATION (WHO) С
 - 1. Guideline Value (GV)
 - 2. Tentative GV
 - Aesthetic GV
- D US ENVIRONMENTAL PROTECTION AGENCY (EPA)
 - 1. Maximum Contaminant Level (MCL)
 - Suggested No-Adverse Effect Level (SNAEL)
 - 3. Lifetime Health Advisory

 - 4. EPA Ambient Water Quality Criteria 47. EPA Ambient Water Quality Criteria for Total PAH
- EUROPEAN ECONOMIC COMMUNITY (EEC)
 - Health Related Guideline Level
 Aesthetic Guideline Level

 - 3. Maximum Admissable Concentration (MADC)
- G CALIFORNIA STATE DEPARTMENT OF HEALTH-GUIDELINE VALUE
- I NEW YORK STATE AMBIENT WATER GUIDELINE
- N/A NONE AVAILABLE

LABORATORY RESULTS, REMARK DESCRIPTIONS

	No Sample Taken
BDL	Below Minimum Measurement Amount
<t< td=""><td>Greater Than Detection Limit But Not Confident (SEE INTERPRETATION OF RESULTS ABOVE)</td></t<>	Greater Than Detection Limit But Not Confident (SEE INTERPRETATION OF RESULTS ABOVE)
>	Results Are Greater Than The Upper Limit
<=>	Approximate Result
!CS	No Data: Contamination Suspected
! IL	No Data: Sample Incorrectly Labelled
! 15	No Data: Insufficient Sample
. VI !	No Data: Inverted Septum
!LA	No Data: Laboratory Accident
! LD	No Data: Test Queued After Sample Discarded
!NA	No Data: No Authorization To Perform Reanalysis
INP	No Data: No Procedure
!NR	No Data: Sample Not Received
! OP	No Data: Obscured Plate
! QU	No Data: Quality Control Unacceptable
!PE	No Data: Procedural Error - Sample Discarded
! PH	No Data: Sample pH Outside Valid Range
!RE	No Data: Received Empty
!RO	No Data: See Attached Report (no numeric results)
! SM	No Data: Sample Missing
! \$\$	No Data: Send Separate Sample Properly Preserved
!UI	No Data: Indeterminant Interference
1TX	No Data: Time Expired
A3C	Approximate, Total Count Exceeded 300 Colonies
APL	Additional Peak, Large, Not Priority Pollutant
APS	Additional Peak, Less Than, Not Priority Pollutant
C1C	Possible Contamination, Improper Cap
CRO	Calculated Result Only
PPS	Test Performed On Preserved Sample
RMP	P and M-Xylene Not Separated
RRV	Rerun Verification
RVU	Reported Value Unusual
SPS	Several Peaks, Small, Not Priority Pollutant

- Unreliable: Could Not Confirm By Reanalysis UCR UCS Unreliable: Contamination Suspected Unreliable: Indeterminate Interference UIN Positive After X Number Of Hours
- XP
- (TO6) Result Taken After # Hours Τ#

DISTRIBUTION SYSTEM

PAU	TREATED
N /N	

FECA	L COLIFORM MF	BACTERIOLOGICAL (CT/100ML)		Det'n limit = 0	GUIDELINE = 0 (A1)
N	ov	0	•		
STAN	DRD PLATE CNT	MF (COUNTS/ML)		DET'N LIMIT = 0	GUIDELINE = 500/ML (A3)
J N	UL IOV	:	10 0		
TOT	L COLIFORM MF	(CT/100ML)		DET'N LIMIT = 0	GUIDELINE = 5/100ML(A1)
h	VO	0	•		
тсо	LIFORM BCKGRD	MF (CT/100ML)		DET'N LIMIT = 0	GUIDELINE = N/A
k	ov	0	•		

RAM	TREATED

	CHEMISTRY (FLD)		
FLD CHLORINE	(COMB) (MG/L)	DET'N LIMIT = 0	GUIDELINE = N/A
JAN		.200		
MAP		200		
MAY	•	100		
PAT	•	. 100		
JUL	•	.900		
SEP		.100		
NOV		1.100		
FLD CHLORINE	FREE (MG/L)		DET'N LIMIT = 0	GUIDELINE = N/A
JAN		.700		
MAR		800		
MAY		700		
111	•	.700		
JUL	•	-000		
SEP	•	.500		
NOV	•	-900		
FLD CHLORINE	(TOTAL) (MG/L)	DET'N LIMIT = 0	GUIDELINE = N/A
JAN		900		
MAD	•	1 000		
MAY	•	1.000		
MAT	•	.000		
JUL	•	1.500		
SEP		.600		
NOV		2.000		
FLD PH (DMNS	LESS)		DET'N LIMIT = N/A	GUIDELINE = 6.5-8.5(A4)
LAN	8 400	7 600		
MAD	8 100	7.000		
MAK	8.100	7.400		
MAY	8.200	7.300		
JUL	8.200	7.400		
SEP	8.500	7.500		
NOV	8.400	7.600		
FLD TEMPERAT	URE (DEG.C)		DET'N LIMIT = N/A	GUIDELINE = 15 (A3)
LAN	500	1 200		
JAN		1.200		
HAK	1.000	2.000		
MAY	9.000	10.500		
JUL	17.000	17.500		
SEP	21,000	21,200		
NOV	11.000	11,200		
FLD TURBIDIT	Y (FTU)		DET'N LIMIT = N/A	GUIDELINE = 1 (A1)
JAN	1.900	.140		
MAR	3,400	-		
MAY	3 600	120		
1111	2 100	600		
000	7.500	.000		
SEP	5.500	.050		
NOV	4.300	.300		
	•••••			

TABLE 5 DRINKING WATER SURVEILLANCE PROGRAM WALPOLE ISLAND WTP 1990

WATER TREATMENT PLANT

RAW

TREATED

	CNEMISTRY	(1 AR)		
ALKALINITY	(MG/L)		DET'N LIMIT = 0.2	GUIDELINE = $30-500$ (A4)
JAN	84.000	77.600		
MAR	91.000	81.000		
MAY	85.400	72.300		
JUL	85.000	79.000		
SEP	83.000	77.500		
NOV	84.100	77.000		
CALCIUM (M	IG/L)		DET'N LIMIT = 0.2	GUIDELINE = 100 (F2)
	70 (00	74 700		
JAN	30.600	31.300		
MAK	35.700	35.800		
MAT	28.800	29.100		
JUL	28.800	30.200		
SEP	29.200	30.300		
NUV	30.000	50.400		
CHLORIDE (MG/L)		DET'N LIMIT = 0.2	GUIDELINE = 250 (A3)
JAN	11,500	12,100		
MAR	10.300	10,800		
MAY	9,100	9,900		
.00	9 100	0 000		
SEP	9 200	10 400		
NOV	6,100	7,000		
COLOUR (HZ	U)		DET'N LIMIT = 0.5	GUIDELINE = 5 (A3)
JAN	1.000 <t< td=""><td>.500 <t< td=""><td></td><td></td></t<></td></t<>	.500 <t< td=""><td></td><td></td></t<>		
MAR	BDL	3.500		
MAY	.500 <t< td=""><td>.500 <t< td=""><td></td><td></td></t<></td></t<>	.500 <t< td=""><td></td><td></td></t<>		
JUL	1.000 <t< td=""><td>BDL</td><td></td><td></td></t<>	BDL		
SEP	.500 <t< td=""><td>BDL</td><td></td><td></td></t<>	BDL		
NON	.500 <t< td=""><td>.500 <t< td=""><td></td><td></td></t<></td></t<>	.500 <t< td=""><td></td><td></td></t<>		
CONDUCTIVI	TY (UMHO/CM)		DET'N LIMIT = 1.	GUIDELINE = 40D (F2)
JAN	235	2/2		
MAR	246	2/2		
MAY	228	274		
111	220	230		
SED	22/	224		
NOV	219	231		
	ARBON (MG/I)			
0.00 ond C			DET N EINT100	GOIDELINE - J.O (AJ)
JAN	1.600	1.200		
MAR	1.300	1.100		
MAY	1.700	1.300		
JUL	1.600	1.400		
SEP	1,600	1,200		
NOV	1,600	1,200		

TABLE 5

DRINKING WATER SURVEILLANCE PROGRAM WALPOLE ISLAND WTP 1990

WATER TREATMENT PLANT

DISTRIBUTION SYSTEM

		RAW	TREATED	
FLUORI	DE (MG/L)		DET'N LIMIT = 0.01	GUIDELINE = 2.4 (A1)
JAN	.080	.080		
MAK	.080	.060		
JUL SEP	.080	.080 .080		
NOV	.060	.100		
HARDNE	SS (MG/L)		DET'N LIMIT = 0.5	GUIDELINE = 80-100 (A4)
JAN	107.600	109.300		
HAY	103.900	104.300		
JUL	105.000	108.000		
NOV	105.000	106.000		
IONCAL	(DMNSLESS)		DET'N LIMIT = N/A	GUIDELINE = N/A
JAN	3.895	3.911		
MAY	.438	1.017		
JUL	.913	2.133		
NOV	1.731 3.140	2.583		
LANGEL	IERS INDEX (DMNSL	.ESS)	DET'N LIMIT = N/A	GUIDELINE = N/A
JAN	.286	.099		
MAR	.314	. 126		
MAY	.180	041		
SEP	. 185	.099		
NOV	.184	.047		
MAGNES	IUM (MG/L)		DET'N LIMIT = 0.1	GUIDELINE = 30 (F2)
JAN	7.550	7.550		
MAR	7.900	7.900		
JUL	8.000	8.000		
SEP	7.700	7.600		
NUV	7.400	7.400		
SODIUM	(MG/L)		DET'N LIMIT = 0.2	GUIDELINE = 200 (A4)
JAN	7.200	7.300		
MAY	5.720	5.600		
JUL	5.800	5.800		
SEP	5.200	5.700		
		4.000		

TABLE 5 DRINKING WATER SURVEILLANCE PROGRAM WALPOLE ISLAND WTP 1990

TREATED

WATER TREATMENT PLANT

DISTRIBUTION SYSTEM

GUIDELINE = 0.05 (F2)

GUIDELINE = 1

(A1)

	u.
2.0	

A

- -

N

MAR

MAY

JUL

SEP NOV

JAN

MAR

MAY

JUL

SEP

NOV

JAN

MAR

MAY

SEP

NOV

JAN

MAR

.

PH (DMNSLESS)

TOTAL NITRATES (MG/L

NITROGEN TOT KJELD (MG/L

.002 <T

.004 <T

.004 <T

.004 <T

.003 <T

.310

.370

.355

.300

.275

.270

.180

.170

.160

.190

.170

.190

8.290

8.220

.........

)

)

MONIUM	TOTAL	(MG/L)		DET'N LIMIT = 0.002
JAN		.008	<1		BDL	
MAR		BDL			BDL	
MAY		.012			BDL	
JUL		.014			BDL	
SEP		.008	<ī		BDL	
NOV		.010			.006 <t< td=""><td></td></t<>	
TRITE	(MG/L)			•••••	DET'N LIMIT = 0.001
JAN		.003	<t< td=""><td></td><td>BDL</td><td></td></t<>		BDL	

.001 <T

.002 <T

.001 <T .002 <T

.

BDL

.310

.355

.310

.290

.295

.110 .090 <T

.130

.150

8.130 8.080

.

.070 <T

DET'N LIMIT = 0.005 GUIDELINE = 10 (A1) DET'N LIMIT = 0.02 GUIDELINE = N/A DET'N LIMIT = N/A GUIDELINE = 6.5-8.5(A4)

MAY 8.200 8.050 JUL 8.280 8.250 8.210 8.140 SEP NOV 8.190 8.090 . DET'N LIMIT = 0.0005 PHOSPHORUS FIL REACT (MG/L) JAN BDL BDL .000 <T .000 <T MAR MAY .001 <T .001 <T JUL BDL BDL SEP BDL BDL .001 <T .004 NOV

GUIDELINE = N/A

RAW TREATED

PHOSPHORUS	TOTAL (MG/L)	DET'N LIMIT = 0.002	GUIDELINE = .40 (F2)
JAN	.007 <t< th=""><th>.002 <1</th><th></th><th></th></t<>	.002 <1		
MAR	-006 <t< th=""><th>RDL</th><th></th><th></th></t<>	RDL		
MAY	.008 <t< th=""><th>.002 <t< th=""><th></th><th></th></t<></th></t<>	.002 <t< th=""><th></th><th></th></t<>		
JUL	.004 <t< th=""><th>BDL</th><th></th><th></th></t<>	BDL		
SEP	.007 <t< th=""><th>.004 <t< th=""><th></th><th></th></t<></th></t<>	.004 <t< th=""><th></th><th></th></t<>		
NOV	.008 <t< th=""><th>BDL</th><th></th><th></th></t<>	BDL		
SULPHATE (MG/L)		DET'N LIMIT = .200	GUIDELINE = 500 (A3)
JAN	16.460	23,590		
MAR	16.290	24.570		
MAY	16.750	28,940		
JUL	16.760	23.310		
SEP	16.270	22,150		
NOV	16.630	26.130		
TURBIDITY	(FTU)		DET'N LIMIT = 0.05	GUIDELINE = 1 (A1)
JAN	1,200	.150 <t< th=""><th></th><th></th></t<>		
MAR	2,900	.540		
MAY	3.400	.200 <t< th=""><th></th><th></th></t<>		
JUL	1.400	.230 <t< th=""><th></th><th></th></t<>		
SEP	2.400	.220		
NOV	2.400	.260		

RAW	TREATED

	METALS		DET'N LIMIT = 0.10	GUIDELINE = 100 (A4)
JAN	13.000	38.000		
MAR	28.000	45.000		
MAY	51.000	39.480		
JUL	23.000	140.000		
SEP	41.000	270.000		
NOV	36.000	120.000		
ARSENIC (L	JG/L)		DET'N LIMIT = 0.10	GUIDELINE = 25 (A1)
JAN	.510 <t< td=""><td>.220 <t< td=""><td></td><td></td></t<></td></t<>	.220 <t< td=""><td></td><td></td></t<>		
MAR	.660 <t< td=""><td>.220 <t< td=""><td></td><td></td></t<></td></t<>	.220 <t< td=""><td></td><td></td></t<>		
MAY	.830 <t< td=""><td>.600 <t< td=""><td></td><td></td></t<></td></t<>	.600 <t< td=""><td></td><td></td></t<>		
JUL	.280 <t< td=""><td>.220 <t< td=""><td></td><td></td></t<></td></t<>	.220 <t< td=""><td></td><td></td></t<>		
SEP	.610 <t< td=""><td>.490 <t< td=""><td></td><td></td></t<></td></t<>	.490 <t< td=""><td></td><td></td></t<>		
NOV	.450 <t< td=""><td>.380 <t< td=""><td></td><td></td></t<></td></t<>	.380 <t< td=""><td></td><td></td></t<>		
BARIUM (UG	5/L ()		DET'N LIMIT = 0.05	GUIDELINE = 1000 (A2)
JAN	15,000	15,000		
MAR	14.000	13,000		
MAY	15,000	13.650		
JUL	15.000	14.000		
SEP	14.000	14.000		
NOV	14.000	13.000		
BORDN (UG/	′L)		DET'N LIMIT = 2.00	GUIDELINE = 5000 (A1)
JAN	15.000 <t< td=""><td>16.000 <t< td=""><td></td><td></td></t<></td></t<>	16.000 <t< td=""><td></td><td></td></t<>		
MAR	16.000 <t< td=""><td>18.000 <t< td=""><td></td><td></td></t<></td></t<>	18.000 <t< td=""><td></td><td></td></t<>		
MAY	43.000	12.650 <t< td=""><td></td><td></td></t<>		
JUL	14.000 <t< td=""><td>16.000 <t< td=""><td></td><td></td></t<></td></t<>	16.000 <t< td=""><td></td><td></td></t<>		
SEP	28,000	28,000		
NOV	15.000 <t< td=""><td>15.000 <t< td=""><td></td><td></td></t<></td></t<>	15.000 <t< td=""><td></td><td></td></t<>		
CADMIUM (L	JG/L)		DET'N LIMIT = 0.05	GUIDELINE = 5 (A1)
JAN	BDL	BDL		
MAR	.060 <t< td=""><td>BDL</td><td></td><td></td></t<>	BDL		
MAY	BDL	.140 <t< td=""><td></td><td></td></t<>		
JUL	BDL	BDL		
SEP	BOI	RDI		
NOV	BDL	BDL		
COBALT (UG	G/L)		DET'N LIMIT = 0.02	GUIDELINE = N/A
JAN	.100 <t< td=""><td>.120 <t< td=""><td></td><td></td></t<></td></t<>	.120 <t< td=""><td></td><td></td></t<>		
MAR	.080 <t< td=""><td>.090 <t< td=""><td></td><td></td></t<></td></t<>	.090 <t< td=""><td></td><td></td></t<>		
MAY	160 <t< td=""><td>280 <t< td=""><td></td><td></td></t<></td></t<>	280 <t< td=""><td></td><td></td></t<>		
.1(1)	150 <t< td=""><td>120 <t< td=""><td></td><td></td></t<></td></t<>	120 <t< td=""><td></td><td></td></t<>		
SEP	060 <t< td=""><td>BDI</td><td></td><td></td></t<>	BDI		
NOV	.070 <t< td=""><td>.030 <t< td=""><td></td><td></td></t<></td></t<>	.030 <t< td=""><td></td><td></td></t<>		

DISTRIBUTION SYSTEM

		RAW	TREATED
CHROMIUM	(UG/L)		DET'N LIMIT = 0.50 GUIDELINE = 50 (A1)
JAN	RDI	BDI	
MAR	BDI	.960	<1
MAY	1.800 <t< th=""><th>BDL</th><th></th></t<>	BDL	
JUL	.510 <t< td=""><td>1.500</td><td><1</td></t<>	1.500	<1
SEP	2.200 <t< td=""><td>2.300</td><td><ī</td></t<>	2.300	<ī
NOV	1.400 <t< td=""><td>1.400</td><td><1</td></t<>	1.400	<1
COPPER (U	G/L)		DET'N LIMIT = 0.50 GUIDELINE = 1000 (A3)
JAN	.890 <t< td=""><td>3.200</td><td><1</td></t<>	3.200	<1
MAR	.900 <t< td=""><td>3.200</td><td><ī</td></t<>	3.200	<ī
MAY	1.900 <t< td=""><td>5.850</td><td></td></t<>	5.850	
JUL	.780 <t< td=""><td>4.400</td><td>ব</td></t<>	4.400	ব
SEP	1.500 <1	1.100	
NOV	i> ۷۷۵.	2.100	<1
IRON (UG/I	L)		DET'N LIMIT = 6.00 GUIDELINE = 300 (A3)
JAN	18.000 <t< b=""></t<>	BDL	
MAR	38.000 <t< td=""><td>BDL</td><td></td></t<>	BDL	
MAY	53.000 <t< td=""><td>BDL</td><td></td></t<>	BDL	
JUL	35.000 <t< td=""><td>BDL</td><td></td></t<>	BDL	
SEP	55.000 <1	BUL 6 100	1
NUV		0.100	
MERCURY (UG/L)		DET'N LIMIT = 0.02 GUIDELINE = 1 (A1)
JAN	.020 <t< td=""><td>.020</td><td><ī</td></t<>	.020	<ī
MAR	BDL	BDL	
MAY	BDL	BDL	
JUL	BDL	BDL	
SEP	BDL	BDL	
NUV	BDL	BUL	
MANGANESE	(UG/L)		DET'N LIMIT = D.05 GUIDELINE = 50 (A3)
JAN	1.400	.670	
MAR	2.300	. 890	
MAY	3.800	.950	
JUL	1.800	.570	
SEP	2.300	.430	<ī
NOV	3.000	.780	
MOLYBDENU	M (UG/L)		DET'N LIMIT = 0.05 GUIDELINE = N/A
JAN	.550	.610	
MAR	.550	.470	<া
MAY	.410 <t< td=""><td>.840</td><td></td></t<>	.840	
JUL	.540	.520	
SEP	.400 <t< td=""><td>.520</td><td></td></t<>	.520	
NOV	.460 <t< td=""><td>.450</td><td></td></t<>	.450	

TABLE 5 DRINKING WATER SURVEILLANCE PROGRAM WALPOLE ISLAND WTP 1990

WATER TREATMENT PLANT

RAW TREATED

NICKEL	(UG/L)		DET'N LIMIT = 0.20	GUIDELINE = 350 (D3)
JAN	1.200 <t< td=""><td>.800 <t< td=""><td></td><td></td></t<></td></t<>	.800 <t< td=""><td></td><td></td></t<>		
MAR	BDL	BDL		
MAY	.260 <t< th=""><th>3.740</th><th></th><th></th></t<>	3.740		
JUL	.330 <t< th=""><th>.620 <t< th=""><th></th><th></th></t<></th></t<>	.620 <t< th=""><th></th><th></th></t<>		
SEP	BDL	BDL		
NOV	ROF	BUL		
LEAD (U	IG/L)		DET'N LIMIT = 0.05	GUIDELINE = 10. (A1)
JAN	.150 <t< td=""><td>.220 <t< td=""><td></td><td></td></t<></td></t<>	.220 <t< td=""><td></td><td></td></t<>		
MAR	.090 <t< td=""><td>.070 <t< td=""><td></td><td></td></t<></td></t<>	.070 <t< td=""><td></td><td></td></t<>		
MAY	.270 <t< td=""><td>.170 <t< td=""><td></td><td></td></t<></td></t<>	.170 <t< td=""><td></td><td></td></t<>		
JUL	.130 <t< td=""><td>.180 <t< td=""><td></td><td></td></t<></td></t<>	.180 <t< td=""><td></td><td></td></t<>		
SEP	.260 <t< td=""><td>.080 <t< td=""><td></td><td></td></t<></td></t<>	.080 <t< td=""><td></td><td></td></t<>		
NOV	.230 <1	.100 <1		
ANTIMON	IY (UG/L)		DET'N LIMIT = 0.05	GUIDELINE = 146 (D4
JAN	.340 <t< td=""><td>.300 <t< td=""><td></td><td></td></t<></td></t<>	.300 <t< td=""><td></td><td></td></t<>		
MAR	.390 <t< td=""><td>.380 <t< td=""><td></td><td></td></t<></td></t<>	.380 <t< td=""><td></td><td></td></t<>		
MAY	.220 <t< td=""><td>.430 <t< td=""><td></td><td></td></t<></td></t<>	.430 <t< td=""><td></td><td></td></t<>		
JUL	.920	.630		
SEP	.460 <t< td=""><td>.460 <1</td><td></td><td></td></t<>	.460 <1		
NUV	.500 <1	.> 00.<1		
SELENIU	JM (UG/L)		DET'N LIMIT = 1.00	GUIDELINE = 10 (A1)
JAN	BDL	BDL		
MAR	BDL	BDL		
MAY	1.200 <t< td=""><td>1.026</td><td></td><td></td></t<>	1.026		
JUL	1.100 <t< td=""><td>1.700 <t< td=""><td></td><td></td></t<></td></t<>	1.700 <t< td=""><td></td><td></td></t<>		
SEP	BUL	BUL		
	DUL	BUL		
STRONT	UM (UG/L)		DET'N LIMIT = 0.10	GUIDELINE = N/A
JAN	100.000	110.000		
MAR	100.000	110.000		
MAY	99.000	105.410		
JUL	100.000	100.000		
SEP	110.000	110.000		
NOV	97.000	97.000		
TITANIU	JM (UG/L)		DET'N LIMIT = 0.50	GUIDELINE = N/A
JAN	2.600 <t< td=""><td>2.200 <t< td=""><td></td><td></td></t<></td></t<>	2.200 <t< td=""><td></td><td></td></t<>		
MAR	5.100	3.900 <t< td=""><td></td><td></td></t<>		
MAY	7.500	7.030		
JUL	4.800 <t< td=""><td>4.300 <t< td=""><td></td><td></td></t<></td></t<>	4.300 <t< td=""><td></td><td></td></t<>		
SEP	5.600	3.10D <t< td=""><td></td><td></td></t<>		
NOV	2.800 <t< td=""><td>2.000 <t< td=""><td></td><td></td></t<></td></t<>	2.000 <t< td=""><td></td><td></td></t<>		

RAW	TREATED

URANIUM (U	G/L)		DET'N LIMIT = 0.05	GUIDELINE = 100 (A1)
JAN	.200 <t< td=""><td>.110 <t< td=""><td></td><td></td></t<></td></t<>	.110 <t< td=""><td></td><td></td></t<>		
MAR	.190 <t< td=""><td>.090 <t< td=""><td></td><td></td></t<></td></t<>	.090 <t< td=""><td></td><td></td></t<>		
MAY	.180 <t< th=""><th>BDL</th><th></th><th></th></t<>	BDL		
JUL	.230 <t< th=""><th>.130 <t< th=""><th></th><th></th></t<></th></t<>	.130 <t< th=""><th></th><th></th></t<>		
SEP	.180 <t< th=""><th>.140 <t< th=""><th></th><th></th></t<></th></t<>	.140 <t< th=""><th></th><th></th></t<>		
NOV	.190 <t< th=""><th>.060 <t< th=""><th></th><th></th></t<></th></t<>	.060 <t< th=""><th></th><th></th></t<>		
VANADIUM (UG/L)		DET'N LIMIT = 0.05	GUIDELINE = N/A
JAN	.270 <t< td=""><td>.170 <t< td=""><td></td><td></td></t<></td></t<>	.170 <t< td=""><td></td><td></td></t<>		
MAR	.260 <t< td=""><td>.170 <t< td=""><td></td><td></td></t<></td></t<>	.170 <t< td=""><td></td><td></td></t<>		
MAY	.270 <t< td=""><td>.140 <t< td=""><td></td><td></td></t<></td></t<>	.140 <t< td=""><td></td><td></td></t<>		
JUL	.250 <t< td=""><td>.160 <t< td=""><td></td><td></td></t<></td></t<>	.160 <t< td=""><td></td><td></td></t<>		
SEP	.310 <t< td=""><td>.190 <t< td=""><td></td><td></td></t<></td></t<>	.190 <t< td=""><td></td><td></td></t<>		
NOV	.190 <t< td=""><td>.140 <t< td=""><td></td><td></td></t<></td></t<>	.140 <t< td=""><td></td><td></td></t<>		
ZINC (UG/L	>		DET'N LIMIT = 0.20	GUIDELINE = 5000 (A3)
JAN	1.300 <t< td=""><td>5.600</td><td></td><td></td></t<>	5.600		
MAR	1.700 <t< td=""><td>2.300</td><td></td><td></td></t<>	2.300		
MAY	3.700	4.750		
JUL	1.500 <t< td=""><td>3.800</td><td></td><td></td></t<>	3.800		
SEP	3.800	45.000		
NOV	2.000 <t< th=""><th>3.000</th><th></th><th></th></t<>	3.000		

DISTRIBUTION SYSTEM

RAW TREATED

		PESTICIDE	S & PCB		
ALPHA BHC	(NG/L)		DET'N LIMIT = 1.000	GUIDELINE = 700 (G
JAN	1.000) <t< td=""><td>1.000 <t< td=""><td></td><td></td></t<></td></t<>	1.000 <t< td=""><td></td><td></td></t<>		
MAR	2.000) <t< td=""><td>1.000 <t< td=""><td></td><td></td></t<></td></t<>	1.000 <t< td=""><td></td><td></td></t<>		
MAY	1.000) <ī	1.000 <t< td=""><td></td><td></td></t<>		
JUL	1.000) <t< td=""><td>2 000 <t< td=""><td></td><td></td></t<></td></t<>	2 000 <t< td=""><td></td><td></td></t<>		
SEP	BD1		BDI		
NON	1.000) <t< td=""><td>1.000 <t< td=""><td></td><td></td></t<></td></t<>	1.000 <t< td=""><td></td><td></td></t<>		

Page 28

RAW TREATED

..... PHENOLICS (UG/L) DET'N LIMIT = .200 GUIDELINE = 2 (A4) IAN BOL BOL

NOV	.800 <t< th=""><th>BDL</th></t<>	BDL
SEP	BDL	BDL
JUL	BDL	BDL
MAY	BDL	BDL
MAR	BDL	.600 <t< th=""></t<>
JAN .	BUL	DUL

DISTRIBUTION SYSTEM

RAW TREATED

	VOLATILES			
RENZENE (UC	VOLATILES		DET'N LIMIT = 0.05	GUIDELINE = 5 (A1)
BENZENE (00				
JAN	.100 <t< td=""><td>IU</td><td></td><td></td></t<>	IU		
MAP	300 <t< td=""><td>.150 <t< td=""><td></td><td></td></t<></td></t<>	.150 <t< td=""><td></td><td></td></t<>		
MAY	BDI	050 <t< td=""><td></td><td></td></t<>		
100	150 <1	100 <t< td=""><td></td><td></td></t<>		
SED	RDI	050 <t< td=""><td></td><td></td></t<>		
NOV	050 <1	BDI		
TOLUENE (UG	/L)		DET'N LIMIT = 0.05	GUIDELINE = 24 (A3)
JAN	.050 <t< td=""><td>!U</td><td></td><td></td></t<>	!U		
MAR	BDL	.050 <t< td=""><td></td><td></td></t<>		
MAY	BDL	BDL		
JUL	.150 <t< td=""><td>.200 <t< td=""><td></td><td></td></t<></td></t<>	.200 <t< td=""><td></td><td></td></t<>		
SEP	BDL	.100 <t< td=""><td></td><td></td></t<>		
NOV	.050 <t< td=""><td>.100 <t< td=""><td></td><td></td></t<></td></t<>	.100 <t< td=""><td></td><td></td></t<>		
ETHYLBENZEN	IE (UG/L)		DET'N LIMIT = 0.05	GUIDELINE = 2.4 (AS)
1411	001			
JAN	BUL	100		
MAK	.050 <1	.100 <1		
MAT	BUL	BUL		
JUL	BDL	.050 <1		
SEP	BOL	.100 <1		
NOV	.050 <t< td=""><td>.05D <t< td=""><td></td><td></td></t<></td></t<>	.05D <t< td=""><td></td><td></td></t<>		
M-XYLENE (U	JG/L)		DET'N LIMIT = 0.10	GUIDELINE = 300 (A3*)
144	BDI	111		
MAR	BDI	BDL		
MAY	BOL	BD1		
.1111	200 <t< td=""><td>.100 <t< td=""><td></td><td></td></t<></td></t<>	.100 <t< td=""><td></td><td></td></t<>		
SEP	BDI	BOI		
NOV	BDL	.100 <t< td=""><td></td><td></td></t<>		
O-XYLENE (U	IG/L)		DET'N LIMIT = 0.05	GUIDELINE = 300 (A3*)
JAN	BDL	in		
MAR	BDL	BDL		
MAY	BDL	BDL		
JUL	.100 <t< td=""><td>.050 <t< td=""><td></td><td></td></t<></td></t<>	.050 <t< td=""><td></td><td></td></t<>		
SEP	BDL	BDL		
NOV	BDL	BDL		
STYRENE (UG	;/L)		DET'N LIMIT = 0.05	GUIDELINE = 100 (D1)
JAN	050 <t< td=""><td>111</td><td></td><td></td></t<>	111		
MAP	100 <t< td=""><td>RDI</td><td></td><td></td></t<>	RDI		
MAY	RDI	BDL		
111	050 <t< td=""><td>BOL</td><td></td><td></td></t<>	BOL		
CED	BDI	BDI		
NOV	100 <1	BDL		
	. 100 - 1			

TABLE 5 DRINKING WATER SURVEILLANCE PROGRAM WALPOLE ISLAND WTP 1990

TREATED

WATER TREATMENT PLANT

RAW

DISTRIBUTION SYSTEM

METHYLENE	CHLORIDE (UG/L)	DET'N LIMIT = 0.50	GUIDELINE = 50 (A1)
JAN	BOL	10		
MAR	BDL	BDL		
MAY	3.000 <t< th=""><th>BDL</th><th></th><th></th></t<>	BDL		
JUL	BDL	BDL		
SEP	BDL	BDL		
NUV	BUL	DVL		
CHLOROFORM	1 (UG/L)		DET'N LIMIT = 0.10	GUIDELINE = 350 (A1+)
JAN	BOL	!U		
MAR	BDL	10.500		
MAY	.200 <t< th=""><th>10.400</th><th></th><th></th></t<>	10.400		
JUL	BDL	19.000		
SEP	BDL	14.400		
NOV	BDL	13.200		
111, TRIC	HLOROETHANE (UG/L)	DET'N LIMIT = 0.02	GUIDELINE = 200 (D1)
JAN	BDL	!U		
MAR	BOL	BDL		
MAY	.100 <t< th=""><th>BDL</th><th></th><th></th></t<>	BDL		
JUL	BDL	BDL		
SEP	BDL	BDL		
NOV	BDL	.020 <t< td=""><td></td><td></td></t<>		
DICHLOROB	ROMOMETHANE (UG/L)	DET'N LIMIT = 0.05	GUIDELINE = 350 (A1+)
JAN	BOL	i U		
MAR	BDL	7.300		
MAY	BDL	10.600		
JUL	BDL	9.250		4
SEP	BDL	8.000		
NOV	BOL	7.000		
CHLOROD I BI	ROMOMETHANE (UG/L)	DET'N LIMIT = 0.10	GUIDELINE = 350 (A1+)
JAN	BDL	!U	·	
MAR	BDL	3.400		
MAY	BDL	8.900		
JUL	BDL	4.000		
SEP	BOL	3.800		
NOV	BDL	3.200		
T-CHLOROE	THYLENE (UG/L))	DET'N LIMIT = 0.05	GUIDELINE = 5 (D1
JAN	.050 <t< td=""><td>!U</td><td></td><td></td></t<>	!U		
MAR	BOL	BDL		
MAY	BOL	BDL		
JUL	BOL	BDL		
SEP	BOL	BDL		
NON	BDL	BDL		

TABLE 5

DRINKING WATER SURVEILLANCE PROGRAM WALPOLE ISLAND WTP 1990

WATER TREATMENT PLANT

RAW TREATED

.

DISTRIBUTION SYSTEM

BROMOFORM (UG/L)		DET'N LIMIT = 0.20	GUIDELINE = 350 (A1+)
JAN	BDL	iU		
MAR	BDL	.400 <t< td=""><td></td><td></td></t<>		
MAY	BDL	1.600 <t< td=""><td></td><td></td></t<>		
JUL	BDL	.400 <t< td=""><td></td><td></td></t<>		
SEP	BDL	.400 <t< td=""><td></td><td></td></t<>		
NOV	BDL	.400 <t< td=""><td></td><td></td></t<>		
TOTL TRIHAL	OMETHANES (UG/L	>	DET'N LIMIT = 0.50	GUIDELINE = 350 (A1)
JAN	BDL	iU		
MAR	BDL	21.600		
MAY	BDL	22,500		
JUL	BDL	32,650		
SEP	BDL	26,550		
NOV	BDL	23.850		

TRACE LEVELS OF TOLUENE ARE LABORATORY ARTIFACTS DERIVED FROM THE ANALYTICAL METHODOLOGY.

TRACE LEVELS OF STYRENE ARE CONSIDERED TO BE LABORATORY ARTIFACTS RESULTING FROM THE LABORATORY SHIPPING CONTAINERS.

		DETECTION	
SCAN/PARAMETER	UNIT	LIMIT	GUIDELINE
		•••••	
0.0000000000000000000000000000000000000			
BACTERIOLOGICAL			
FECAL COLLEODY NEWDOWNE STLEDATION	CT (100M)	0	0 (41)
CTANDADD DLATE COUNT MEMODANE FILT	CT/HU	0	500/HL (A3)
TOTAL COLLEODY PACKCOCIND NS	CT/100MI	0	SUU/ML (AS)
TOTAL COLLEGON MENDRAVE ELLIDATION	CT/100HL	ő	5/100ML (A1)
TOTAL COLITORA HEADRARE FILTRATION	CT/ TOORL	Ŭ	J/ 100AL (A1)
CHEMISTRY (FLD)			
FIELD COMBINED CHLORINE RESIDUAL	MG/L	0	N/A
FIELD TOTAL CHLORINE RESIDUAL	MG/L	0	N/A
FIELD FREE CHLORINE RESIDUAL	MG/L	0	N/A
FIELD PH	DMNSLESS	N/A	6.5-8.5 (A3)
FIELD TEMPERATURE	DEG.C	N/A	15.0 (A3)
FIELD TURBIDITY	FTU	N/A	1.0 (A1)
CHEWISTRY (LAR)			
ALKALINITY	MG/L	0.2 3	30-500 (A3)
AMHONIUM TOTAL	MG/L	0.002	0.05 (F2)
CALCIUM	MG/L	0.2	100 (F2)
CHLORIDE	MG/L	0.2	250 (A3)
COLOUR	TCU	0.5	5.0 (A3)
CONDUCTIVITY	UMHO/CM	1.0	400 (F2)
CYANIDE COOLUGE CLOBON	MG/L	0.001	U.2 (A1)
DISSOLVED UKGANIC CARBON	MG/L	0.1	5.0 (AS)
NARDNESS	MG/L	0.01	2.4 (AI) 20-100 (A()
IANCELLEDS INDEX	MG/L	0.5 0	N/A
MAGNESTIM	MG/I	0 1	30 0 (F2)
NITRITE	MG/L	0 001	1.0 (A1)
NITROGEN TOTAL KJELDAHL	MG/L	0.02	N/A
PH	DMNSLESS	N/A	6.5-8.5 (A4)
PHOSPHORUS FIL REACT	MG/L	0.0005	N/A
PHOSPHORUS TOTAL	MG/L	0.002	0.4 (F2)
SODIUM	MG/L	0.2	200 (A4)
SULPHATE	MG/L	0.2	500 (A3)
TOTAL NITRATES	MG/L	0.005	10.0 (A1)
TURBIDITY	FTU	0.05	1.0 (A1)
CHLOROAROMATICS			
CHEOROAROPATIES			
123 TRICHLOROBENZENE	NG/L	5.0	N/A
1234 TETRACHLOROBENZENE	NG/L	1.0	N/A
1235 TETRACHLOROBENZENE	NG/L	1.0	N/A
124 TRICHLOROBENZENE	NG/L	5.0	10000 (I)
1245-TETRACHLOROBENZENE	NG/L	1.0	38000 (D4)
135 TRICHLOROBENZENE	NG/L	5.0	N/A
236 TRICHLOROTOLUENE	NG/L	5.0	N/A
245 TRICHLOROTOLUENE	NG/L	5.0	N/A
ZOA TRICHLORUTULUENE	NG/L	5.0	10 (01)
	NG/L	1.0	(50 (0()
	NG/L	5.0	206000 (04)
NEXACILOROGICEOPERTADIERE	NG/L	1.0	1900 (04)
DCTACHLOROSTYRENE	NG/L	1.0	N/A
PENTACHLOROBENZENE	NG/L	1.0	74000 (D4)
UNLOKUPHENULS			
234 TRICHLOROPHENOL	NG/L	100.0	N/A
2345 TETRACHLOROPHENOL	NG/L	20.0	N/A
2356 TETRACHLOROPHENOL	NG/L	10.0	N/A

SCAN / DADANETER	UNIT	LIMIT	GUIDELINE
245 TRICHLOROPHENOL	NG/L	100.0	2600000 (D4)
246 TRICHLOROPHENOL	NG/L	20.0	5000 (A1)
PENTACHLOROPHENOL	NG/L	10.0	60000 (A1)
METALS			
ALUHINUK	UG/L	0.10	100 (A4)
ANTIHONY	UG/L	0.05	146 (D4)
ARSENIC	UG/L	0.10	25 (A1) 1000 (A2)
BERYLLIN	UG/L	0.05	6800 (D4)
BORON	UG/L	2.00	5000 (A1)
CADMIUM	UG/L	0.05	5 (A1)
CHROMIUM	UG/L	0.50	50 (A1)
CORDER		0.02	1000 (43)
IPON		6.00	300 (A3)
LEAD	UG/L	0.05	10 (A1)
MANGANESE	UG/L	0.05	50 (A3)
MERCURY	UG/L	0.02	1 (A1)
HOLTBDENOM	UG/L	0.05	350 (03)
SELENIUM	UG/L	1.00	10 (A1)
SILVER	UG/L	0.05	50 (A1)
STRONTIUM	UG/L	0.10	N/A
THALLIUM	UG/L	0.05	13 (D4)
	UG/L	0.50	N/A 100 (A1)
VANADIUM	UG/L	0.05	N/A
ZINC	UG/L	0.20	5000 (A3)
PAH			
ANTNDACENE	NG/I	1.0	N/A
BENZO(A) ANTHRACENE	NG/L	20.0	N/A
BENZO(A) PYRENE	NG/L	5.0	10.0 (A1)
BENZO(B) CHRYSENE	NG/L	2.0	N/A
BENZO(B) FLUORANTHENE	NG/L	10.0	N/A ·
BENZO(C) PIRENE	NG/L NG/L	20.0	N/A
BENZO(K) FLUORANTHENE	NG/L	1.0	N/A
CHRYSENE	NG/L	50.0	N/A
CORONENE	NG/L	10.0	N/A
DIBENZU(A,H) ANTHRAGENE	NG/L	10.0	N/A
FLUORANTHENE	NG/L	20.0	42000,0 (D4)
INDENO(1,2,3-C,D) PYRENE	NG/L	20.0	N/A
PERYLENE	NG/L	10.0	N/A
PHENANTHRENE	NG/L	10.0	N/A
	NG/L	20.0	8/A
FLOTICIDES & FUB			
ALACHLOR (LASSO)	NG/L	500.0	5000 (A2)
	NG/L	1.0	700 (A1) 700 (C)
ALPHA REXACTLOROCICLOREXANE (BRC)	NG/L NG/I	2.0	7000 (41)
AMETRINE	NG/L	50.0	300000 (D3)
ATRATONE	NG/L	50.0	N/A
ATRAZINE	NG/L	50.0	60000 (A2)
DES ETHYL ATRAZINE	NG/L	200.0	60000 (A2)
CYANAZINE (BLADEX)	NG/L	100.0	10000 (A2)
O,P-DDD	NG/L	5.0	10 (1)
DIELDRIN	NG/L	2.0	700 (A1)
ENDOSULFAN 1 (THIODAN I)	NG/L	2.0	74000 (D4)
ENDOSULFAN 2 (THIODAN II)	NG/L	5.0	74000 (04)

4

		DETECTION	
SCAN/PARAMETER	UNIT	LIMIT	GUIDELINE
ENDOSIN FAN SUNPHATE (THIODAN SUNPHATE)	NG/I	5.0	N/A
ENDDIN	NG/L	5.0	1600 (03)
CANNA CHI ODDANE	NC /I	2.0	7000 (41)
GARRA CREOKDARE	NG/L	2.0	7000 (41)
HEPTACHLOK	NG/L	1.0	3000 (A1)
REPTACHLOR EPOXIDE	NG/L	1.0	3000 (A1)
LINDANE (GAMMA BHC)	NG/L	1.0	4000 (A1)
METHOXYCHLOR	NG/L	5.0	900000 (A1)
METOLACHLOR	NG/L	500.0	50000 (A2)
METRIBUZIN (SENCOR)	NG/L	100.0	80000 (A1)
MIREX	NG/L	5.0	N/A
P,P-DDD	NG/L	5.0	N/A
O.P-DDT	NG/L	5.0	30000 (A1)
DXYCHLORDANE	NG/L	2.0	N/A
PCB	NG/I	20.0	3000 (A2)
PPDDE	NG/L	1.0	30000 (41)
PPDDT	NG/L	5.0	30000 (41)
PROMETONE	NG/L	50.0	52500 (03)
DOMETRANE	NG/L	50.0	1000 (42)
PROPERTINE	NG/L	50.0	700000 (07)
PROFAZINE	NG/L	50.0	10000 (03)
SIMAZINE	NG/L	50.0	10000 (A2)
D-ETHYL SIMAZINE	NG/L	200.0	10000 (A2)
TOXAPHENE	NG/L	500.0	5000 (A1)
PHENOLICS			
PHENOLICS (UNFILTERED REACTIVE)	UG/L	0.2	2 (A4)
SPECIFIC PESTICIDES			
2.4 D PROPIONIC ACID	NG/I	100.	N/A
2 4 5-TRICHLOROPHENOXY ACETIC ACID	NG/L	50.	280000 (41)
2 4-DICHIOPORITYPIC ACID (2 4-D)	NG/L	100	100000 (41)
24-DICHLORODDHENOYYDUTYDIC ACID (24-DD)	NG/L	200	18000 (83)
DITVIATE (DITAN)	NC /I	2000	2/5000 (03)
CARRADYL (CEVIN)	NG/L	2000.	240000 (00)
CARDACIL (SEVIR)	NG/L	200.	90000 (A1)
CARBUFURAN	NG/L	2000.	90000 (AT)
CREORPTRIFUS (DURSBAN)	NG/L	20.	N/A
CICP (CHLORPROPHAM)	NG/L	2000.	350000 (6)
UIALLATE	NG/L	2000.	N/A
DIAZINON	NG/L	20.	20000 (A1)
DICAMBA	NG/L	50.	120000 (A1)
DICHLOROVOS	NG/L	20.	N/A
EPTAM	NG/L	2000.	N/A
ETHION	NG/L	20.	35000 (G)
IPC	NG/L	2000.	N/A
MALATHION	NG/L	20.	190000 (A1)
METHYL PARATHION	NG/L	50.	7000 (83)
METRYLTRITHION	NG/L	20.	N/A
MEVINPHOS	NG/I	20.	N/A
PAPATHION	NG/I	20.	50000 (A1)
DHODATE (THINET)	NG/L	20	2000 (42)
PRODOVID (PAYCON)	NC/L	2000	140000 (03)
PEDAN (BAIGON)	NG/L	2000.	N/A
DOWNEL	NG/L	20.	N/A
CILVER () (E TD)	NG/L	20.	10000 (41)
SILVEX (2,4,5-1P)	NG/L	20.	10000 (AT)
VOLATILES			
1,1 DICHLOROETHANE	UG/L	0.10	N/A
1,1 DICHLOROETHYLENE	UG/L	0.10	7 (01)
1,2 DICKLOROBENZENE	UG/L	0.05	200 (A1)
1.2 DICHLOROETHANE	UG/L	0.05	5 (A1)

.

		DETECTION	
SCAN/PARAMETER	UNIT	LIMIT	GUIDELINE
1,2 DICHLOROPROPANE	UG/L	0.05	5 (D1)
1,3 DICHLOROBENZENE	UG/L	0.10	3750 (05)
1,4 DICHLOROBENZENE	UG/L	0.10	5 (A1)
111, TRICHLOROETHANE	UG/L	0.02	200 (D1)
112 TRICHLOROETHANE	UG/L	0.05	0.6 (04)
1122 TETRACHLOROETHANE	UG/L	0.05	0.17(D4)
BENZENE	UG/L	0.05	5 (A1)
BROMOFORM	UG/L	0.20	350 (A1+)
CARBON TETRACHLORIDE	UG/L	0.20	5 (A1)
CHLOROBENZENE	UG/L	0.10	1510 (D3)
CHLORODIBROMOMETHANE	UG/L	0.10	350 (A1+)
CHLOROFORM	UG/L	0.10	350 (A1+)
DICHLOROBROMOMETHANE	UG/L	0.05	350 (A1+)
ETHLYENE DIBROMIDE	UG/L	0.05	50 (D1)
ETHYLBENZENE	UG/L	0.05	2.4 (A3)
M-XYLENE	UG/L	0.10	300 (A3*)
METHYLENE CHLORIDE	UG/L	0.50	50 (A1)
O-XYLENE	UG/L	0.05	300 (A3*)
P-XYLENE	UG/L	0.10	300 (A3*)
STYRENE	UG/L	0.05	100 (D1)
TETRACHLOROETHYLENE	UG/L	0.05	5 (D1)
TRANS 1,2 DICHLOROETHYLENE	UG/L	0.10	70 (D1)
TOLUENE	UG/L	0.05	24 (A3)
TOTAL TRIHALOMETHANES	UG/L	0.50	350 (A1)
TRICHLOROETHYLENE	UG/L	0.10	50 (A1)
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DRINKING WATER SURVEILLANCE PROGRAM PROGRAM DESCRIPTION

The Drinking Water Surveillance Program (DWSP) for Ontario monitors drinking water quality at municipal water supply systems. The DWSP Database Management System provides a computerized drinking water quality information system for the supplies monitored. The objectives of the program are to provide:

- immediate, reliable, current information on drinking water quality;
- a flagging mechanism for guideline exceedance;
- a definition of contaminant levels and trends;
- a comprehensive background for remedial action;
- a framework for assessment of new contaminants; and
- an indication of treatment efficiency of plant processes.

PROGRAM

The DWSP officially began in April 1986 and is designed to eventually include all municipal water supplies in Ontario. In 1990, 76 systems were being monitored. Water supply locations have been prioritized for surveillance based primarily on criteria such as population density, probability of contamination and geographical location.

An ongoing assessment of future monitoring requirements at each location will be made. Monitoring will continue at the initial locations at an appropriate level and further locations will be phased into the program as resources permit.

A major goal of the program is to collect valid water quality data in context with plant operational characteristics at the time of sampling. As soon as sufficient data have been accumulated and analyzed, both the frequency of sampling and the range of parameters may be adjusted accordingly.

Assessments are carried out at all locations prior to initial sampling, in order to acquire complete plant process and distribution system details and to designate (and retrofit if necessary) all sampling systems and locations. This ensures that the sampled water is a reflection of the water itself. Samples are taken of raw (ambient water) and treated water at the treatment plant and of consumer's tap water in the distribution system. In order to determine possible effects of distribution on water quality, both standing and free flow water in old and new sections of the distribution system are sampled. Sampling is carried out by operational personnel who have been trained in applicable procedures.

Comprehensive standardized procedures and field test kits are supplied to sampling personnel. This ensures that samples are taken and handled according to standard protocols and that field testing will supply reliable data. All field and laboratory analyses are carried out using "approved documented procedures". Most laboratory analyses are carried out by the Ministry of Environment (MOE), Laboratory Services Branch. Radionuclides are analyzed by the Ministry of Labour.

DATA REPORTING MECHANISM

When the analytical results are transferred from the MOE laboratory into the DWSP system, printouts of the completed analyses are sent to the MOE District Officer, the appropriate operational staff and are also retained by the DWSP unit.

PROGRAM INPUTS AND OUTPUTS

There are four major inputs and four major outputs in the program.

Program Input - Plant and Distribution System Description

The system description includes plant specific non-analytical information acquired through a questionnaire and an initial plant visit. During the initial assessment of the plant and distribution system, questionnaire content is verified and missing information added. It is intended that all data be kept current with scheduled annual updates.

The Plant and Distribution System Description consists of the following seven components:

1. PROCESS COMPONENT INVENTORY

All physical and chemical processes to which the water is subjected, from the intake pipe to the consumers' tap (where possible), are documented. These include: process type, general description of physical structures, material types, sizes, and retention time for each process within the plant. The processes may be as simple as transmission or as complex as carbon adsorption.

2. TREATMENT CHEMICALS

Chemicals used in the treatment processes, their function, application point, supplier and brand-name are recorded. Chemical dosages applied on the day of sampling are recorded in DWSP.

3. PROCESS CONTROL MEASUREMENTS

Documentation of in-plant monitoring of process parameters (eg. turbidity, chlorine residuals, pH, aluminum residuals) including methods used, monitoring locations and frequency is contained in this section. Except for the recorded Field Data, in-plant monitoring results are not retained in DWSP but are retained by the water treatment plant personnel.

4. DESIGN FLOW AND RETENTION TIME

Hydraulic capacity, designed and actual, is noted here. Retention time (the time that a block of water is retained in the plant) is also noted. Maximum, minimum and average flow, as well as a record of the flow rate on the day of sampling, are recorded in DWSP.

5. DISTRIBUTION SYSTEM DESCRIPTION

This area includes the storage and transmission characteristics of the distribution system after the water leaves the plant.

6. SAMPLING SYSTEM

Each plant is assessed for its adequacy in terms of the sampling of bacteriological, organic and inorganic parameters. Prime considerations in the assessment and design of the sampling system are:

- i/ the sample is an accurate representation of the actual water condition, eg. raw water has had no chemical treatment;
- ii/ the water being sampled is not being modified by the sampling system;
- iii/ the sample tap must be in a clean area of the plant, preferably a lab area; and
 - iv/ the sample lines must be organically inert (no plastic, ideally stainless steel).

It is imperative that the sampled water be a reflection not of the sampling system but of the water itself.

The sampling system documentation includes: origin of the water; date sampling was initiated; size, length and material type (intake,

discharge and tap); pump characteristics (model, type, capacity); and flow rate.

7. PERSONNEL

This section contains the names, addresses and phone numbers of current plant management and operational staff, distribution system management and operational staff, Medical Officer of Health and appropriate MOE personnel associated with the plant.

Program Input - Field Data

The second major input to DWSP is field data. Field data is collected at the plant and from the distribution system sites on the day of sampling. Field data consists of general operating conditions and the results of testing for field parameters. General operating conditions include chemicals used, dosages, flow and retention time on the day of sampling, as well as, monthly maximum, minimum and average flows. Field parameters include turbidity, chlorine residuals (free, combined and total), temperature and pH. These parameters are analyzed according to standardized DWSP protocols to allow for interplant comparison.

Program Input - Laboratory Analytical Data

The third major input to DWSP is Laboratory Analytical Data. Samples gathered from the raw, treated and distribution sampling sites are analyzed for the presence of approximately 180 parameters at a frequency of two to twelve times per year. Sixty-five percent of the parameters are organic. Parameters measured may have health or aesthetic implications when present in drinking water. Many of the parameters may be used in the treatment process or may be treatment by-products. Due to the nature of certain analytical instruments, parameters may be measured in a "scan" producing some results for parameters that are not on the DWSP priority list, but which may be of interest. The majority of parameters are measured on a routine basis. Those that are technically more difficult and/or costly to analyze, however, are done less frequently. These include Specific Pesticides and Chlorophenols.

Although the parameter list is extensive, additional parameters with the potential to cause health or aesthetic related problems may be added provided reliable analytical and sampling methods exist.

All laboratory generated data is derived from standardized, documented analytical protocols. The analytical method is an integral part of the data and as methods change, notation will be made and comparison data documented.

Program Input - Parameter Reference Information

The fourth major input to DWSP is Parameter Reference Information. This is a catalogue of information for each substance analyzed on DWSP. It includes parameter name and aliases, physical and chemical properties, basic toxicology, world-wide health limits, treatment methods and uses. The Parameter Reference Information is computerized and can be accessed through the Query function of the DWSP database. An example is shown in figure 1.

Program output - Query

All DWSP information is easily accessed through the Query function, therefore, anything from addresses of plant personnel to complete water quality information for a plant's water supply is instantly available. The DWSP computer system makes relatively complex inquiries manageable. A personal password allowing access into the DWSP query mode in all MOE offices is being developed by the DWSP group.

Program Output - Action Alerts

Drinking Water quality in Ontario is evaluated against provincial objectives as outlined in the Ontario Drinking Water Objectives publication. Should the reported level of a substance in treated water exceed the Ontario Drinking Water Objective, an "Action Alert" requiring resampling and confirmation is issued. This assures that operational staff, health authorities and the public are notified as soon as possible of the confirmation of an exceedance and remedial action taken. This report supplies a history of the occurrence of past exceedances at the plant plus a historical summary on the parameter of concern.

In the absence of Ontario Drinking Water Objectives, guidelines/limits from other agencies are used. The Parameter Listing System, published by MOE (ISBN 0-7729-4461-X), catalogues and keeps current guidelines for 650 parameters from agencies throughout the world. If these guidelines are exceeded, the results are flagged and evaluated by DWSP personnel. An "Action Alert" will be issued if warranted.

Program Output - Report Generation

Custom reports can be generated from DWSP to meet MOE Regional needs and to respond to public requests.

Program Output - Annual Reports

It is the practice of DWSP to produce an annual report containing analytical data along with companion plant information.

FIG.1

MOE - DRINKING WATER ASSESSMENT PROGRAM (DWSP)

PARAMETER REFERENCE INFORMATION

BENZENE (B2001P)

VOLATILES

0111001	************		1001 100000	01111			
SOURCE	FROM	то	METHOD	GUIDELINE	UNIT	NOTE	
CAL C	85/01			0.700	µg/L	AL	
CDWG C	87/01			5.000	µg/L	MAC	
EPA C	87/07			5.000	µg/L	MCL	
EPAA C	80/11			6.600	µg/L	AMBIENT *	ŧ *
FERC C	84/05			1.000	µg/L	MCL	
WHO C	84/01			10.000	µg/L	GV	

DESCRIPTION: NAME: BENZENE

CAS#: 71-43-2

MOLECULAR FORMULAE: C6H6

CLASS HEALTH METHOD DOCODO LINTT UG/L

DETECTION LIMIT: (FOR METHOD POCODO) 0.05 µg/L

- SYNONYMS: BENZOL; BENZOLE; COAL NAPHTHA; CARBON OIL (27). CYCLOHEXATRIENE (41).
- CHARACTERISTICS: COLOURLESS TO LIGHT-YELLOW, MOBILE, NON-POLAR LIQUID, OF HIGHLY REFRACTIVE NATURE, AROMATIC ODOUR; VAPOURS BURN WITH SMOKING FLAME (30).

PROPERTIES: SOLUBILITY IN WATER: 1780-1800 mg/L AT 25C (41).
THRESHOLD ODOUR: 0.5 - 10 PPM IN WATERTHRESHOLD TASTE:
0.5 mg/L IN WATER (39).
ENVIRONMENTAL FATE: MAY BIOACCUMULATE IN LIVING
ORGANISMS AND APPEARS TO ACCUMULATE IN ANIMAL

TISSUES THAT EXHIBIT A HIGH LIPID CONTENT OR REPRESENT MAJOR METABOLIC SITES, SUCH AS LIVER OR BRAIN; SMALL QUANTITIES EVAPORATE FROM SOILS OR ARE DEGRADED RATHER QUICKLY (80).

SOURCES: COMMERCIAL: PETROLEUM REFINING; SOLVENT RECOVERY; COAL TAR DISTILLATION (39); FOOD PROCESSING AND TANNING INDUSTRIES; COMBUSTION OF CAR EXHAUST. ENVIRONMENTAL: POSSIBLE SOURCE IS RUNOFF.

- USES: DETERGENTS; NYLON; INTERMEDIATE IN PRODUCTION OF OTHER COMPOUNDS, SUCH AS PESTICIDES; SOLVENT FOR EXTRACTION AND RECTIFICATION IN RUBBER INDUSTRY; DEGREASING AND CLEANSING AGENT; GASOLINE.
- TOXICITY: RATING: 4 (VERY TOXIC).

ACUTE: IRRITATING TO MUCOUS MEMBRANES; SYMPTOMS INCLUDE RESTLESSNESS, CONVULSIONS, EXCITEMENT, DEPRESSION; DEATH MAY FOLLOW RESPIRATORY FAILURE. CHRONIC: MAY CAUSE ANAEMIA AND LEUKAEMIA (45); MUTAGENIC. MODE OF ACTION: CHROMOABERRATION IN LYMPHOCYTE

CULTURES.

CARCINOGENICITY: A KNOWN HUMAN CARCINOGEN.

REMOVAL: THE FOLLOWING PROCESSES HAVE BEEN SUCCESSFUL IN REMOVING BENZENE FROM WASTEWATER: GAC ADSORPTION, PRECIPITATION WITH ALUM AND SUBSEQUENT REMOVAL VIA SEDIMENTATION, COAGULATION AND FLOCCULATION, SOLVENT EXTRACTION, OXIDATION

ADDITIONAL PROPERTIES:

MOLECULAR WEIGHT: 78.12 MELTING POINT: 5.5°C (27). BOILING POINT: 80.1°C (27). SPECIFIC GRAVITY: 0.8790 AT 20°C (27). VAPOUR PRESSURE: 100 MM AT 26.1°C (27). HENRY'S LAW CONSTANT: 0.00555 ATM-M3/MOLE (41). LOG OCT./WATER PARTITION COEFFICIENT: 1.95 TO 2.13 (39). CARBON ADSORPTION: K=1.0; 1/N=1.6; R=0.97; PH=5.3 (41) SEDIMENT/WATER PARTITION COEFFICIENT: NO DATA

NOTES: EPA PRIORITY POLLUTANT.

DWSP SAMPLING GUIDELINE

i) Raw and Treated at Plant

General Chemistry	-500 mL plastic bottle (PET 500) -rinse bottle and cap with sample water three times -fill to 2 cm from top
Bacteriological	 -220 mL plastic bottle with white seal on cap -do <u>not</u> rinse bottle, preservative has been added -avoid touching bottle neck or inside of cap -fill to top of red label as marked
Metals	-500 mL plastic bottle (PET 500) -rinse bottle and cap three times -fill to 2 cm from top -add 10 drops nitric acid (HNO ₃) (Caution: HNO ₃ is corrosive)
Volatiles (duplicates) (OPOPUP)	 -45 mL glass vial with septum (teflon side must be in contact with sample) -do <u>not</u> rinse bottle -fill bottle completely without bubbles
Organics (OWOC),(OWTRI),(OAPAHX)	<pre>-1 L amber glass bottle per scan -do <u>not</u> rinse bottle -fill to 2 cm from top -when 'special pesticides' are requested three extra bottles must be filled</pre>

Cyanide	-500 mL plastic bottle (PET 500) -rinse bottle and cap three times -fill to 2 cm from top -add 10 drops sodium hydroxide (NaOH) (Caution: NaOH is corrosive)
Mercury	-250 mL glass bottle -rinse bottle and cap three times -fill to top of label -add 20 drops each nitric acid (HNO ₃) and potassium dichromate (K ₂ Cr ₂ O ₇) (Caution: HNO ₃ &K ₂ Cr ₂ O ₇ are corrosive)
Phenols	-250 mL glass bottle -do <u>not</u> rinse bottle, preservative has been added -fill to top of label
Radionuclides (as scheduled)	-4 L plastic jug -do <u>not</u> rinse, carrier added -fill to 5 cm from top
Organic Characterization (GC/MS - once per year)	 -1 L amber glass bottle; instructions as per organic -250 mL glass bottle -do not rinse bottle -fill completely without bubbles

Steps:

- 1. Let sampling water tap run for an adequate time to clear the sample line.
- 2. Record time of day on submission sheet.
- 3. Record temperature on submission sheet.
- 4. Fill up all bottles as per instructions.
- Record chlorine residuals (free, combined and total for treated water only), turbidity and pH on submission sheet.

ii) Distribution Samples (standing water)

General	Chemistry	-500 mL plastic bottle (PET 500) -rinse bottle and cap with sample water three times -fill to 2 cm from top
Metals		-500 mL plastic bottle (PET 500) -rinse bottle and cap three times -fill to 2 cm from top -add 10 drops nitric acid (HNO ₃) (Caution: HNO ₃ is corrosive)

Steps:

- 1. Record time of day on submission sheet.
- 2. Place bucket under tap and open cold water.
- 3. Fill to predetermined volume.
- 4. After mixing the water, record the temperature on the submission sheet.
- 5. Fill general chemistry and metals bottles.
- Record chlorine residuals (free, combined and total), turbidity and pH on submission sheet.

iii) Distribution Samples (free flow)

General Chemistry	-500 mL plastic bottle (PET 500) -rinse bottle and cap with sample water three times -fill to 2 cm from top
Bacteriological	-250 mL plastic bottle with white seal on cap -do <u>not</u> rinse bottle, preservative has been added -avoid touching bottle neck or inside of cap -fill to top of red label as marked

Page 47

Metals	-500 mL plastic bottle (PET 500) -rinse bottle and cap three times -fill to 2 cm from top -add 10 drops nitric acid HNO ₃ (Caution: HNO ₃ is corrosive)
Volatiles (duplicate) (OPOPUP)	 -45 mL glass vial with septum (teflon side must be in contact with sample) -do <u>not</u> rinse bottle, preservative has been added -fill bottle completely without bubbles
Organics (OWOC) (OAPAHX)	-1 L amber glass bottle per scan -do <u>not</u> rinse bottle -fill to 2 cm from top

Steps:

- 1. Record time of day on submission sheet.
- 2. Let cold water flow for five minutes.
- 3. Record temperature on submission sheet.
- 4. Fill all bottles as per instructions.
- 5. Record chlorine residuals (free, combined and total), turbidity and pH on submission sheet.